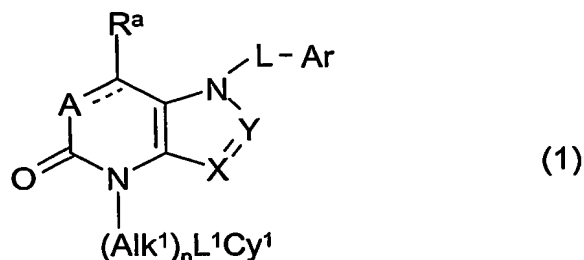


**Claims:**

1. A compound of formula (1):



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wherein:

the dashed line joining A and C(R<sup>a</sup>) is present and represents a bond and A is a -N= atom or a -C(R<sup>b</sup>)= group, or the dashed line is absent and A is a -C(R<sup>b</sup>)(R<sup>c</sup>)- or -N(R<sup>d</sup>)- group;

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R<sup>a</sup>, R<sup>b</sup> and R<sup>c</sup> is each independently a hydrogen or halogen atom or an optionally substituted alkyl, -CN, -CO<sub>2</sub>R<sup>1</sup> (where R<sup>1</sup> is a hydrogen atom or an optionally substituted alkyl group) or -CONR<sup>1</sup>R<sup>2</sup> group (where R<sup>2</sup> is a hydrogen atom or an optionally substituted alkyl group);

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R<sup>d</sup> is a hydrogen atom or an alkyl group;

X and Y is each a nitrogen atom or a -C(R<sup>e</sup>)= or -C(Alk<sup>2</sup>R<sup>e</sup>)= group [where Alk<sup>2</sup> is an optionally substituted alkylene, alkenylene or alkynylene chain and R<sup>e</sup> is a hydrogen or halogen atom or a -CN, -OR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -C(X<sup>a</sup>)R<sup>1</sup> (where X<sup>a</sup> is an oxygen or sulphur atom), -Cy<sup>2</sup> (where Cy<sup>2</sup> is an optionally substituted, saturated or unsaturated non-aromatic carbocyclic ring optionally containing one or more -O-, -S-, -NH- or -C(X<sup>a</sup>)- atoms or groups), -NR<sup>1a</sup>R<sup>2a</sup> (where R<sup>1a</sup> and R<sup>2a</sup>, which may be the same or different, is each a hydrogen atom or an optionally substituted alkyl or Cy<sup>2</sup> group, or together with the nitrogen atom to which they are attached form an optionally substituted, saturated or unsaturated cyclicamino ring optionally containing one or more -O- or -S- atoms or -NH- or -C(X<sup>a</sup>)- groups), -C(X<sup>a</sup>)NR<sup>1a</sup>R<sup>2a</sup>, -S(O)<sub>2</sub>NR<sup>1a</sup>R<sup>2a</sup>, -N(R<sup>3a</sup>)C(X<sup>a</sup>)R<sup>1</sup> (where R<sup>3a</sup> is a hydrogen atom or an optionally substituted alkyl group), -N(R<sup>3a</sup>)C(X<sup>a</sup>)NR<sup>1a</sup>R<sup>2a</sup>, -N(R<sup>3a</sup>)S(O)<sub>2</sub>R<sup>1</sup>, -N[S(O)<sub>2</sub>R<sup>1</sup>]<sub>2</sub>, -N(R<sup>3a</sup>)S(O)<sub>2</sub>NR<sup>1a</sup>R<sup>2a</sup>, -N(R<sup>3a</sup>)C(O)OR<sup>1</sup>, -N(R<sup>3a</sup>)C(NR<sup>1</sup>)NR<sup>1a</sup>R<sup>2a</sup>, -C(R<sup>1</sup>)NOR<sup>2</sup>, -C(NR<sup>1</sup>)NR<sup>1a</sup>R<sup>2a</sup>, -C(X<sup>a</sup>)NR<sup>1a</sup>OR<sup>2a</sup> or -C(O)N(R<sup>3a</sup>)NR<sup>1a</sup>R<sup>2a</sup> group];

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L is a -C(O)-, -C(S)- or -C(R<sup>1f</sup>)(R<sup>1g</sup>)- group (where R<sup>1f</sup> and R<sup>1g</sup>, which may be the same or different, is each a hydrogen atom or a straight or branched C<sub>1-3</sub>alkyl group optionally substituted by one, two or three fluorine atoms, or R<sup>1f</sup> and R<sup>1g</sup> together with the carbon atom to which they are attached form a cyclopropyl group), or a -CH<sub>2</sub>CH<sub>2</sub>- group;

5 n is zero or the integer 1;

Alk<sup>1</sup> is an optionally substituted aliphatic or heteroaliphatic chain;

L<sup>1</sup> is a covalent bond or a linker atom or group;

Cy<sup>1</sup> is an optionally substituted cycloaliphatic, polycycloaliphatic, heterocycloaliphatic, polyheterocycloaliphatic, aromatic or heteroaromatic group, or is  
10 additionally a hydrogen atom when n is the integer 1 and/or L<sup>1</sup> is a linker atom or group;  
and

Ar is an optionally substituted aromatic or heteroaromatic group;

provided that the compound of formula (1) is other than 3,7-dibenzyl-3,7-dihydro-2H-purinone;

15 and the salts, solvates, hydrates and N-oxides thereof.

2. A compound as claimed in claim 1 wherein the dashed line joining A and C(R<sup>a</sup>) is present and represents a bond and A is a -C(R<sup>b</sup>)- group, in which R<sup>a</sup> and R<sup>b</sup> are as defined in claim 1.

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3. A compound as claimed in claim 2 wherein R<sup>a</sup> and R<sup>b</sup> are both hydrogen.

4. A compound as claimed in any one of the preceding claims wherein X is a -CH= group.

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5. A compound as claimed in any one of the preceding claims wherein Y is a -C(R<sup>e</sup>)= group, in which R<sup>e</sup> is hydrogen, -CN, -COR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -CONR<sup>1a</sup>R<sup>2a</sup>, -S(O)<sub>2</sub>NR<sup>1a</sup>R<sup>2a</sup>, -CONR<sup>1a</sup>OR<sup>2a</sup> or -C(O)N(R<sup>3a</sup>)NR<sup>1a</sup>R<sup>2a</sup>, and R<sup>1</sup>, R<sup>1a</sup>, R<sup>2a</sup> and R<sup>3a</sup> are as defined in claim 1.

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6. A compound as claimed in any one of the preceding claims wherein Cy<sup>1</sup> is phenyl, methylphenyl, methoxyphenyl, thienyl or indolyl.

7. A compound as claimed in any one of the preceding claims wherein Ar represents phenyl, fluorophenyl, difluorophenyl, chlorophenyl, dichlorophenyl, (chloro)(fluoro)phenyl, cyanophenyl, methylphenyl, (fluoro)(methyl)phenyl, methoxyphenyl, nitrophenyl, pyridinyl, chlorothienyl or benzothienyl.

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8. A compound as claimed in claim 1 as herein specifically disclosed in any one of the Examples.

9. A pharmaceutical composition comprising a compound of formula (1) as  
10 defined in claim 1, or a pharmaceutically acceptable salt, solvate, hydrate or *N*-oxide thereof, in association with a pharmaceutically acceptable carrier.

10. The use of a compound of formula (1) as defined in claim 1, or a  
pharmaceutically acceptable salt, solvate, hydrate or *N*-oxide thereof, for the manufacture  
15 of a medicament for the treatment and/or prevention of a disorder for which an inhibitor of p38 kinase is indicated.

11. A method for the treatment and/or prevention of a disorder for which an  
inhibitor of p38 kinase is indicated, which comprises administering to a patient in need of  
20 such treatment a compound of formula (1) as defined in claim 1, or a pharmaceutically acceptable salt, solvate, hydrate or *N*-oxide thereof.